

*Ohio's Full #1 - Claim 19. + Inventor Search*

Kwon 10/694,641

02/12/2005

FILE 'HCAPLUS' ENTERED AT 15:26:10 ON 02 DEC 2005  
L1 75361 S HYPERTENSION+ALL OR HYPERTENSIVE AGENTS+ALL  
L2 33 S L1 AND ?EPOXIDE?(W)?HYDROLASE?

FILE 'REGISTRY' ENTERED AT 15:27:22 ON 02 DEC 2005  
L3 STR  
L4 50 S L3  
L5 99950 S L3 FUL  
L6 STR L3  
L7 50 S L6  
L8 STR L6  
L9 50 S L8

FILE 'HCAPLUS' ENTERED AT 15:35:05 ON 02 DEC 2005  
L10 SELECT RN L2 1-33  
2 S L5 AND L2

FILE 'HCAPLUS' ENTERED AT 15:38:23 ON 02 DEC 2005  
FILE 'REGISTRY' ENTERED AT 15:38:26 ON 02 DEC 2005  
L11 6521 S L8 FUL

FILE 'HCAPLUS' ENTERED AT 15:38:51 ON 02 DEC 2005  
L12 1 S L2 AND L11

FILE 'REGISTRY' ENTERED AT 15:39:38 ON 02 DEC 2005  
L13 STR  
L14 50 S L13  
L15 STR L13  
L16 STR L15  
L17 STR L16  
L18 50 S L17  
DELETE SELECT

FILE 'HCAPLUS' ENTERED AT 15:48:06 ON 02 DEC 2005  
L19 1 S L2 AND (PRD<19980208 OR PD<19980208)  
E KROETZ DEANNA L/AU  
L20 36 S E1-5  
E ZELDIN DARRYL C/AU  
L21 83 S E1-3  
E HAMMOCK BRUCE D/AU  
L22 634 S E1-4  
E MORISSEAU CHRISTOPHE/AU  
L23 58 S E2-5  
L24 2 S L20 AND L21 AND L22 AND L23  
SELECT RN L24 1-2

FILE 'REGISTRY' ENTERED AT 15:50:54 ON 02 DEC 2005  
L25 162 S E1-162

FILE 'HCAPLUS' ENTERED AT 15:52:26 ON 02 DEC 2005  
L26 2 S L24 AND L25  
L27 ANALYZE L26 1-2 CT : 18 TERMS  
L28 6 S L1 AND L11  
L29 4 S L28 AND (PRD<19980208 OR PD<19980208)  
L30 4 S L12 OR L29

FILE 'REGISTRY' ENTERED AT 16:06:19 ON 02 DEC 2005

*I did your  
search in 2  
stages, Thus  
2 search histories*

*Claim 19  
search*

*Inventor  
Search*

*— captured on transcript — results for ②  
(claim 19)  
see attached  
done stat for claim 19*

L31 STR L17  
L32 50 S L31

FILE 'HCAPLUS' ENTERED AT 16:08:12 ON 02 DEC 2005

FILE 'REGISTRY' ENTERED AT 16:10:07 ON 02 DEC 2005

L33 STR L31  
L34 50 S L33  
L35 STR L33  
L36 50 S L35

SAV L35 KWO641L35/L

*saved str. for ①*

FILE 'HCAPLUS' ENTERED AT 16:20:55 ON 02 DEC 2005

SAV L30 KWO641L30/A

=> <-----Page Break----->

=> log hold

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

24.66

1659.09

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

-2.92

-62.78

CA SUBSCRIBER PRICE

*-1300*  
*459*

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 16:22:04 ON 02 DEC 2005

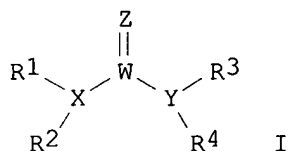
=&gt; d ibib abs hitstr 143 1-2

L43 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2005 ACS on STN  
ACCESSION NUMBER: 2003:197917 HCAPLUS  
DOCUMENT NUMBER: 138:231736  
TITLE: Inhibitors of epoxide hydrolases for the treatment of hypertension  
INVENTOR(S): Kroetz, Deanna L.; Zeldin, Darryl C.  
; Hammock, Bruce D.; Morisseau, Christophe  
PATENT ASSIGNEE(S): Regents of the University of California, USA  
SOURCE: U.S., 36 pp., Cont.-in-part of U.S. 6,150,415.  
CODEN: USXXAM  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 3  
PATENT INFORMATION:

5/4/  
475, 54  
551, 64  
613,  
611, 576,  
588, 585

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6531506	B1	20030311	US 2000-721261	20001121
US 5955496	A	19990921	US 1997-909523	19970812
US 6150415	A	20001121	US 1999-252148	19990218
US 2003119900	A1	20030626	US 2002-328495	20021223
US 6693130	B2	20040217		
US 2004092487	A1	20040513	US 2003-694641	20031027
PRIORITY APPLN. INFO.:			US 1996-23397P	P 19960813
			US 1997-909523	A2 19970812
			US 1999-252148	A2 19990218
			US 2000-721261	A1 20001121
			US 2002-328495	A1 20021223

OTHER SOURCE(S): MARPAT 138:231736  
GI

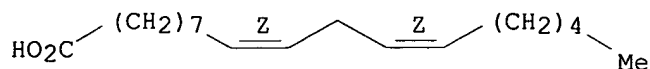


AB The invention provides compds. that inhibit epoxide hydrolase in therapeutic applications for the treatment of hypertension. A preferred class of compds. for practicing the invention have the structure shown by Formula [(R1)(R2)XW(Z)Y(R3)(R4)], wherein Z is oxygen or sulfur, W is carbon phosphorous or sulfur, X and Y is each independently nitrogen, oxygen, or sulfur, and X can further be carbon, at least one of R1-R4 is hydrogen, R2 is hydrogen when X is nitrogen but is not present when X is sulfur or oxygen, R4 is hydrogen when Y is nitrogen but is not present when Y is sulfur or oxygen, R1 and R3 is each independently C1-C20 substituted or unsubstituted alkyl, cycloalkyl, aryl, acyl, or heterocyclic.

IT 60-33-3, Linoleic acid, biological studies 9048-63-9,  
Epoxide hydrolase 97717-69-6, Epoxyeicosatrienoic acid  
192461-94-2 192461-95-3 192461-96-4  
RL: BSU (Biological study, unclassified); BIOL (Biological study)  
(epoxide hydrolases inhibitors for treatment of hypertension)  
RN 60-33-3 HCAPLUS

CN 9,12-Octadecadienoic acid (9Z,12Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 9048-63-9 HCAPLUS

CN Hydratase, epoxide (9CI) (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 97717-69-6 HCAPLUS

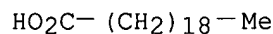
CN Eicosatrienoic acid, epoxy- (9CI) (CA INDEX NAME)

CM 1

CRN 97717-68-5

CMF C20 H38 O3

CCI IDS

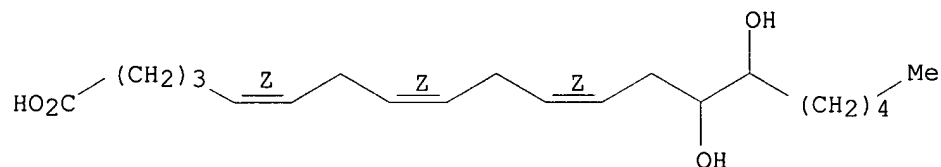


D1-O-D1

RN 192461-94-2 HCAPLUS

CN 5,8,11-Eicosatrienoic acid, 14,15-dihydroxy-, (5Z,8Z,11Z)- (9CI) (CA INDEX NAME)

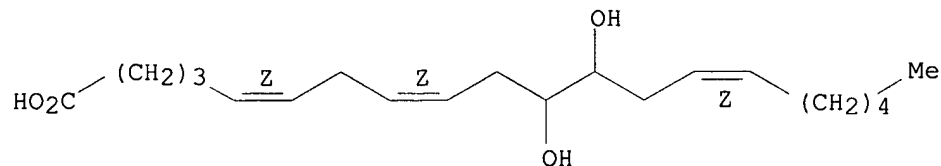
Double bond geometry as shown.



RN 192461-95-3 HCAPLUS

CN 5,8,14-Eicosatrienoic acid, 11,12-dihydroxy-, (5Z,8Z,14Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

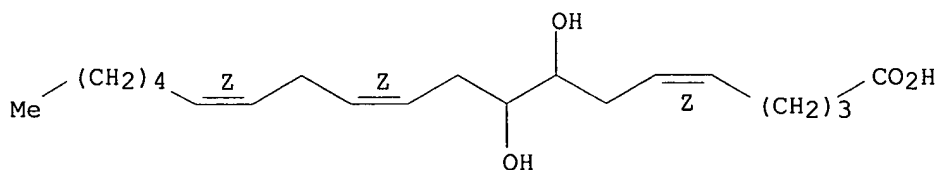


RN 192461-96-4 HCAPLUS

CN 5,11,14-Eicosatrienoic acid, 8,9-dihydroxy-, (5Z,11Z,14Z)- (9CI) (CA INDEX NAME)

INDEX NAME)

Double bond geometry as shown.



IT 101-20-2 538-75-0 886-59-9 1142-07-0  
 1145-53-5 1220-01-5 1885-07-0  
 2222-58-4 2387-23-7 3218-00-6  
 4300-43-0 7570-86-7 13074-28-7  
 13256-78-5 17427-95-1 17661-70-0  
 20258-07-5 25855-24-7 27729-98-2  
 29559-44-2 31510-11-9 34583-53-4  
 36598-10-4 37141-32-5 37141-34-7  
 38652-23-2 53120-54-0 58490-15-6  
 60392-17-8 60392-18-9 62972-70-7  
 70243-17-3 70622-91-2 72802-45-0  
 72802-46-1 75461-45-9 80251-18-9  
 82389-34-2 82499-65-8 89609-46-1  
 91990-71-5 92373-98-3 98819-68-2  
 101452-64-6 104196-23-8 108195-84-2  
 109246-29-9 111681-31-3 118389-04-1  
 120615-85-2 124949-24-2 128276-09-5  
 131519-56-7 148806-83-1 148806-87-5  
 158072-24-3 176327-83-6 179539-07-2  
 195451-88-8 196791-88-5 197357-33-8  
 197455-83-7 200058-85-1 200059-00-3  
 201014-78-0 204922-97-4 246165-79-7  
 272449-13-5 287185-28-8 287185-29-9  
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 338797-34-5 339170-08-0 392713-22-3  
 400839-30-7 401598-13-8 402939-18-8  
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501004-16-6 501004-17-7 501004-18-8

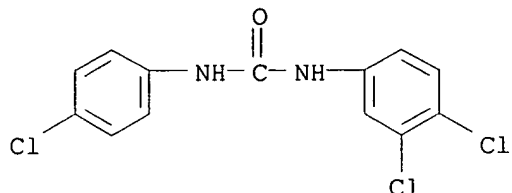
501004-19-9 501004-20-2 501011-71-8

RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL  
(Biological study); USES (Uses)

(epoxide hydrolases inhibitors for treatment of hypertension)

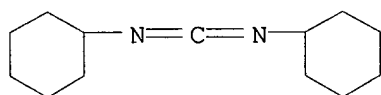
RN 101-20-2 HCAPLUS

CN Urea, N-(4-chlorophenyl)-N'-(3,4-dichlorophenyl)- (9CI) (CA INDEX NAME)



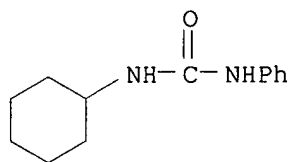
RN 538-75-0 HCAPLUS

CN Cyclohexanamine, N,N'-methanetetraylbis- (9CI) (CA INDEX NAME)



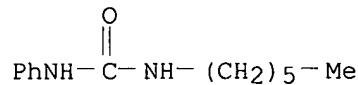
RN 886-59-9 HCAPLUS

CN Urea, N-cyclohexyl-N'-phenyl- (9CI) (CA INDEX NAME)



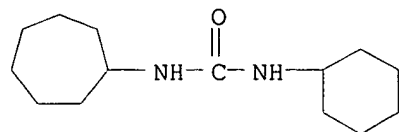
RN 1142-07-0 HCAPLUS

CN Urea, N-hexyl-N'-phenyl- (9CI) (CA INDEX NAME)

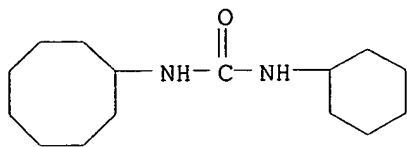


RN 1145-53-5 HCAPLUS

CN Urea, N-cycloheptyl-N'-cyclohexyl- (9CI) (CA INDEX NAME)

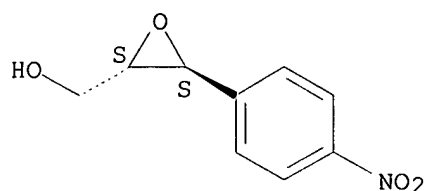


RN 1220-01-5 HCAPLUS  
CN Urea, N-cyclohexyl-N'-cyclooctyl- (9CI) (CA INDEX NAME)

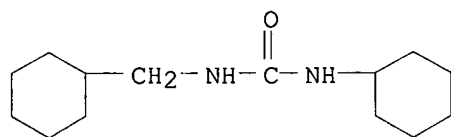


RN 1885-07-0 HCAPLUS  
CN Oxiranemethanol, 3-(4-nitrophenyl)-, (2S,3S)- (9CI) (CA INDEX NAME)

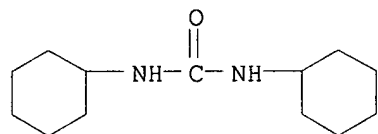
Absolute stereochemistry.



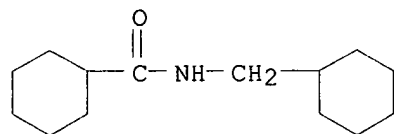
RN 2222-58-4 HCAPLUS  
CN Urea, N-cyclohexyl-N'-(cyclohexylmethyl)- (9CI) (CA INDEX NAME)



RN 2387-23-7 HCAPLUS  
CN Urea, N,N'-dicyclohexyl- (9CI) (CA INDEX NAME)

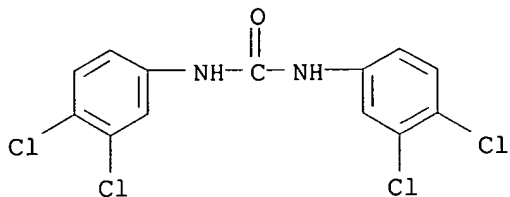


RN 3218-00-6 HCAPLUS  
CN Cyclohexanecarboxamide, N-(cyclohexylmethyl)- (7CI, 8CI, 9CI) (CA INDEX NAME)



RN 4300-43-0 HCAPLUS

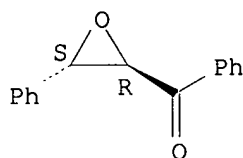
CN Urea, N,N'-bis(3,4-dichlorophenyl)- (9CI) (CA INDEX NAME)



RN 7570-86-7 HCAPLUS

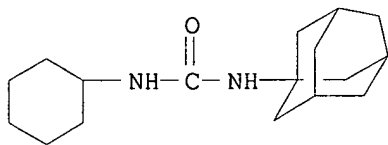
CN Methanone, phenyl[(2R,3S)-3-phenyloxiranyl]-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



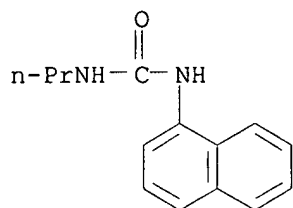
RN 13074-28-7 HCAPLUS

CN Urea, N-cyclohexyl-N'-tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl- (9CI) (CA INDEX NAME)



RN 13256-78-5 HCAPLUS

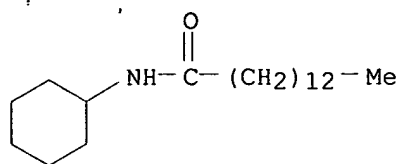
CN Urea, N-1-naphthalenyl-N'-propyl- (9CI) (CA INDEX NAME)



RN 17427-95-1 HCAPLUS

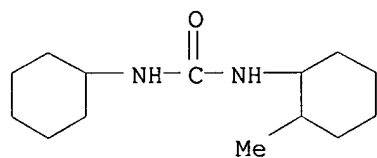
CN Tetradecanamide, N-cyclohexyl- (9CI) (CA INDEX NAME)





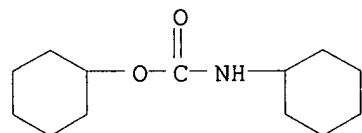
RN 17661-70-0 HCAPLUS

CN Urea, N-cyclohexyl-N'-(2-methylcyclohexyl)- (9CI) (CA INDEX NAME)



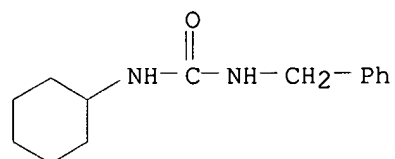
RN 20258-07-5 HCAPLUS

CN Carbamic acid, cyclohexyl-, cyclohexyl ester (9CI) (CA INDEX NAME)



RN 25855-24-7 HCAPLUS

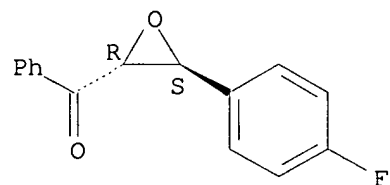
CN Urea, N-cyclohexyl-N'-(phenylmethyl)- (9CI) (CA INDEX NAME)



RN 27729-98-2 HCAPLUS

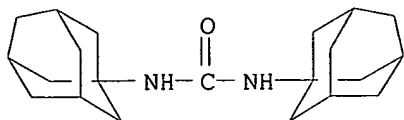
CN Methanone, [(2R,3S)-3-(4-fluorophenyl)oxiranyl]phenyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



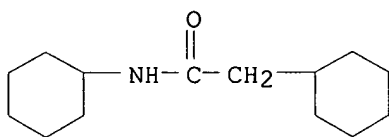
RN 29559-44-2 HCAPLUS

CN Urea, N,N'-bis(tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl)- (9CI) (CA INDEX NAME)



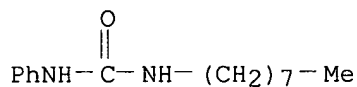
RN 31510-11-9 HCAPLUS

CN Cyclohexaneacetamide, N-cyclohexyl- (8CI, 9CI) (CA INDEX NAME)



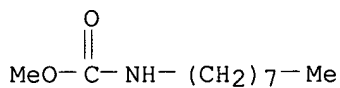
RN 34583-53-4 HCAPLUS

CN Urea, N-octyl-N'-phenyl- (9CI) (CA INDEX NAME)



RN 36598-10-4 HCAPLUS

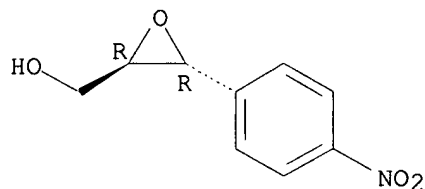
CN Carbamic acid, octyl-, methyl ester (9CI) (CA INDEX NAME)



RN 37141-32-5 HCAPLUS

CN Oxiranemethanol, 3-(4-nitrophenyl)-, (2R,3R)- (9CI) (CA INDEX NAME)

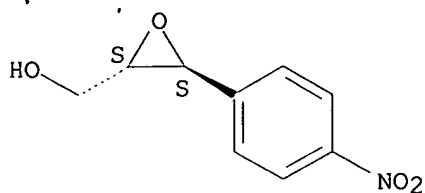
Absolute stereochemistry.



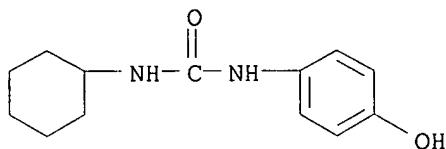
RN 37141-34-7 HCAPLUS

CN Oxiranemethanol, 3-(4-nitrophenyl)-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

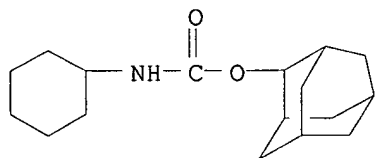
Relative stereochemistry.



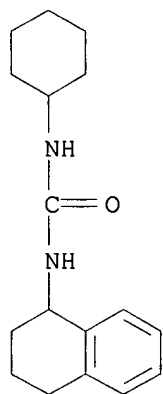
RN 38652-23-2 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-hydroxyphenyl)- (9CI) (CA INDEX NAME)



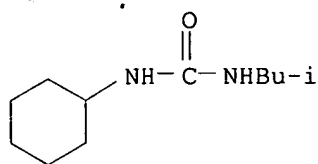
RN 53120-54-0 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, tricyclo[3.3.1.1<sup>3,7</sup>]dec-2-yl ester (9CI) (CA INDEX NAME)



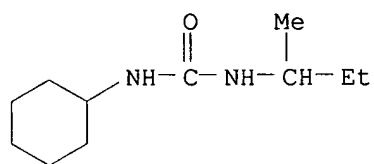
RN 58490-15-6 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(1,2,3,4-tetrahydro-1-naphthalenyl)- (9CI) (CA INDEX NAME)



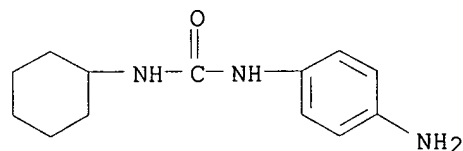
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 CN Urea, N-cyclohexyl-N'-(2-methylpropyl)- (9CI) (CA INDEX NAME)



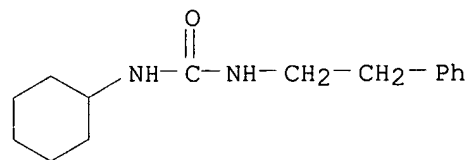
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 CN Urea, N-cyclohexyl-N'-(1-methylpropyl)- (9CI) (CA INDEX NAME)



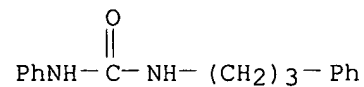
RN 62972-70-7 HCAPLUS  
 CN Urea, N-(4-aminophenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



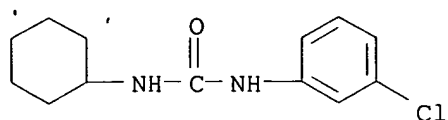
RN 70243-17-3 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(2-phenylethyl)- (9CI) (CA INDEX NAME)



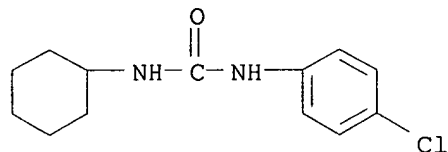
RN 70622-91-2 HCAPLUS  
 CN Urea, N-phenyl-N'-(3-phenylpropyl)- (9CI) (CA INDEX NAME)



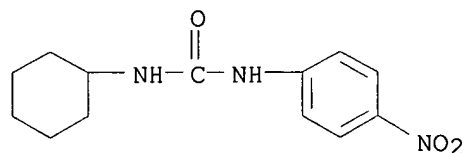
RN 72802-45-0 HCAPLUS  
 CN Urea, N-(3-chlorophenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



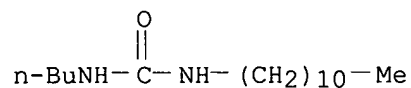
RN 72802-46-1 HCAPLUS  
 CN Urea, N-(4-chlorophenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



RN 75461-45-9 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-nitrophenyl)- (9CI) (CA INDEX NAME)

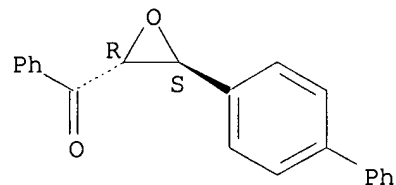


RN 80251-18-9 HCAPLUS  
 CN Urea, N-butyl-N'-undecyl- (9CI) (CA INDEX NAME)

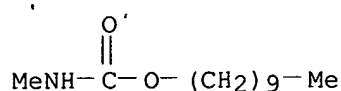


RN 82389-34-2 HCAPLUS  
 CN Methanone, [(2R,3S)-3-[1,1'-biphenyl]-4-yloxiranyl]phenyl-, rel- (9CI)  
 (CA INDEX NAME)

Relative stereochemistry.

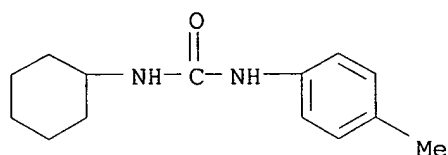


RN 82499-65-8 HCAPLUS  
 CN Carbamic acid, methyl-, decyl ester (9CI) (CA INDEX NAME)



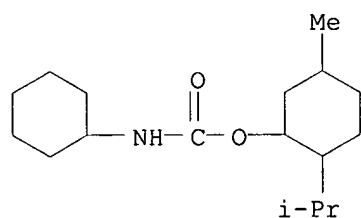
RN 89609-46-1 HCAPLUS

CN Urea, N-cyclohexyl-N'-(4-methylphenyl)- (9CI) (CA INDEX NAME)



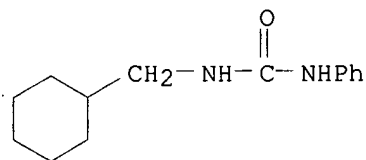
RN 91990-71-5 HCAPLUS

CN Carbamic acid, cyclohexyl-, 5-methyl-2-(1-methylethyl)cyclohexyl ester (9CI) (CA INDEX NAME)



RN 92373-98-3 HCAPLUS

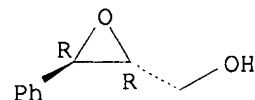
CN Urea, N-(cyclohexylmethyl)-N'-phenyl- (9CI) (CA INDEX NAME)



RN 98819-68-2 HCAPLUS

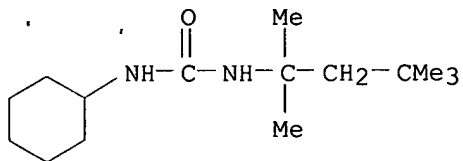
CN Oxiranemethanol, 3-phenyl-, (2R,3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



RN 101452-64-6 HCAPLUS

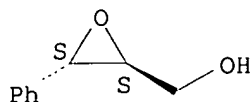
CN Urea, N-cyclohexyl-N'-(1,1,3,3-tetramethylbutyl)- (9CI) (CA INDEX NAME)



RN 104196-23-8 HCAPLUS

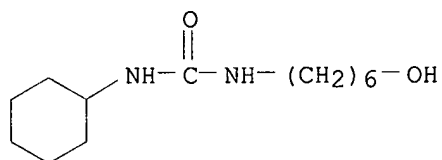
CN Oxiranemethanol, 3-phenyl-, (2S,3S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



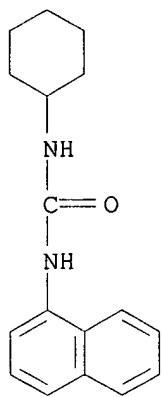
RN 108195-84-2 HCAPLUS

CN Urea, N-cyclohexyl-N'-(6-hydroxyhexyl)- (9CI) (CA INDEX NAME)



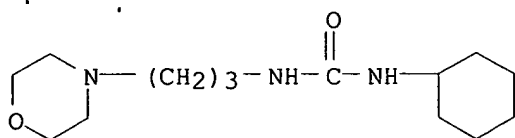
RN 109246-29-9 HCAPLUS

CN Urea, N-cyclohexyl-N'-1-naphthalenyl- (9CI) (CA INDEX NAME)

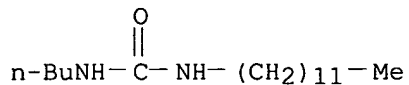


RN 111681-31-3 HCAPLUS

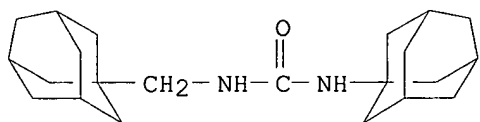
CN Urea, N-cyclohexyl-N'-[3-(4-morpholinyl)propyl]- (9CI) (CA INDEX NAME)



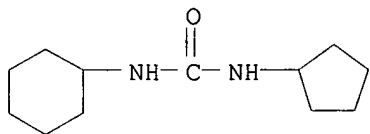
RN 118389-04-1 HCAPLUS  
CN Urea, N-butyl-N'-dodecyl- (9CI) (CA INDEX NAME)



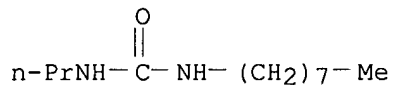
RN 120615-85-2 HCAPLUS  
CN Urea, N-tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl-N'-(tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-ylmethyl)- (9CI) (CA INDEX NAME)



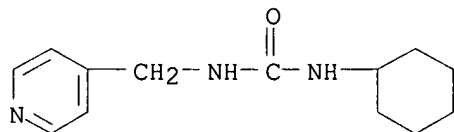
RN 124949-24-2 HCAPLUS  
CN Urea, N-cyclohexyl-N'-cyclopentyl- (9CI) (CA INDEX NAME)



RN 128276-09-5 HCAPLUS  
CN Urea, N-octyl-N'-propyl- (9CI) (CA INDEX NAME)



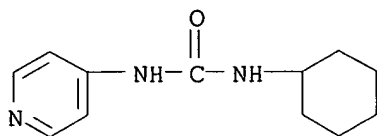
RN 131519-56-7 HCAPLUS  
CN Urea, N-cyclohexyl-N'-(4-pyridinylmethyl)- (9CI) (CA INDEX NAME)





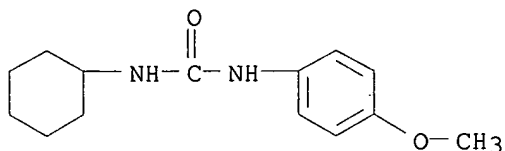
RN 148806-83-1 HCAPLUS

CN Urea, N-cyclohexyl-N'-4-pyridinyl- (9CI) (CA INDEX NAME)



RN 148806-87-5 HCAPLUS

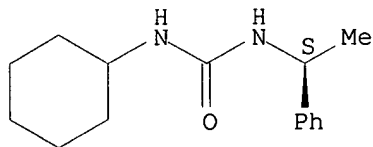
CN Urea, N-cyclohexyl-N'-(4-methoxyphenyl)- (9CI) (CA INDEX NAME)



RN 158072-24-3 HCAPLUS

CN Urea, N-cyclohexyl-N'-[(1S)-1-phenylethyl]- (9CI) (CA INDEX NAME)

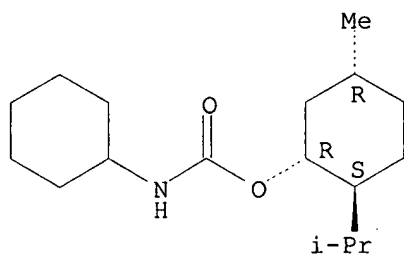
Absolute stereochemistry. Rotation (-).



RN 176327-83-6 HCAPLUS

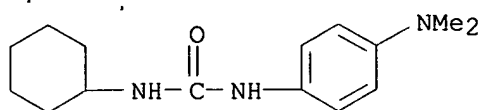
CN Carbamic acid, cyclohexyl-, (1R,2S,5R)-5-methyl-2-(1-methylethyl)cyclohexyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.



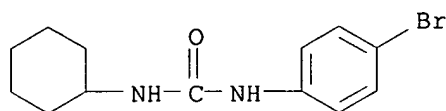
RN 179539-07-2 HCAPLUS

CN Urea, N-cyclohexyl-N'-[4-(dimethylamino)phenyl]- (9CI) (CA INDEX NAME)



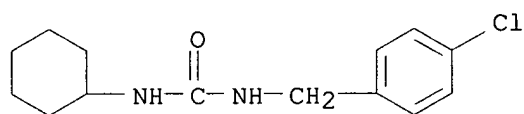
RN 195451-88-8 HCAPLUS

CN Urea, N-(4-bromophenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



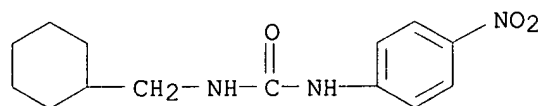
RN 196791-88-5 HCAPLUS

CN Urea, N-[(4-chlorophenyl)methyl]-N'-cyclohexyl- (9CI) (CA INDEX NAME)



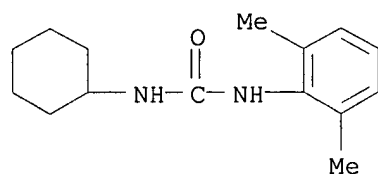
RN 197357-33-8 HCAPLUS

CN Urea, N-(cyclohexylmethyl)-N'-(4-nitrophenyl)- (9CI) (CA INDEX NAME)



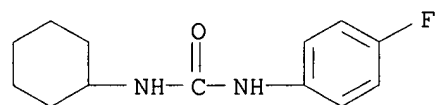
RN 197455-83-7 HCAPLUS

CN Urea, N-cyclohexyl-N'-(2,6-dimethylphenyl)- (9CI) (CA INDEX NAME)



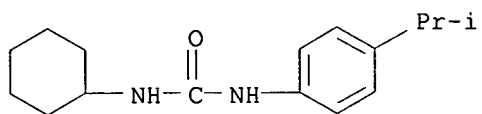
RN 200058-85-1 HCAPLUS

CN Urea, N-cyclohexyl-N'-(4-fluorophenyl)- (9CI) (CA INDEX NAME)



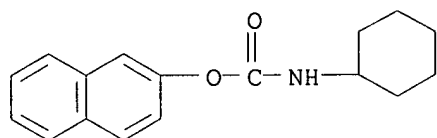
RN 200059-00-3 HCAPLUS

CN Urea, N-cyclohexyl-N'-[4-(1-methylethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 201014-78-0 HCAPLUS

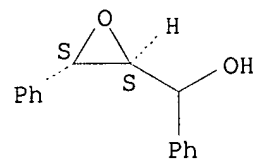
CN Carbamic acid, cyclohexyl-, 2-naphthalenyl ester (9CI) (CA INDEX NAME)



RN 204922-97-4 HCAPLUS

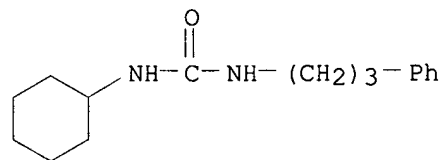
CN Oxiranemethanol,  $\alpha$ ,3-diphenyl-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



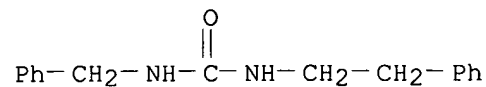
RN 246165-79-7 HCAPLUS

CN Urea, N-cyclohexyl-N'-(3-phenylpropyl)- (9CI) (CA INDEX NAME)



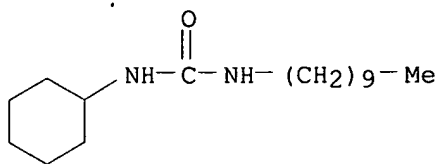
RN 272449-13-5 HCAPLUS

CN Urea, N-(2-phenylethyl)-N'-(phenylmethyl)- (9CI) (CA INDEX NAME)

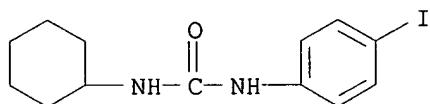


RN 287185-28-8 HCAPLUS

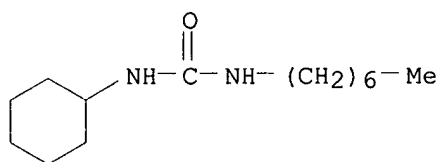
CN Urea, N-cyclohexyl-N'-decyl- (9CI) (CA INDEX NAME)



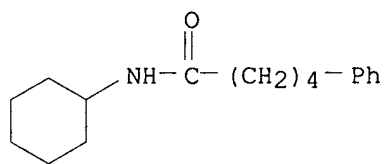
RN 287185-29-9 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-iodophenyl)- (9CI) (CA INDEX NAME)



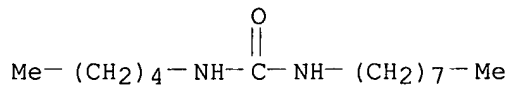
RN 303092-06-0 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-heptyl- (9CI) (CA INDEX NAME)



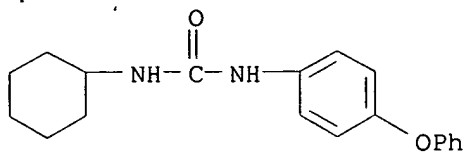
RN 306746-73-6 HCAPLUS  
 CN Benzenepentanamide, N-cyclohexyl- (9CI) (CA INDEX NAME)



RN 306770-63-8 HCAPLUS  
 CN Urea, N-octyl-N'-pentyl- (9CI) (CA INDEX NAME)

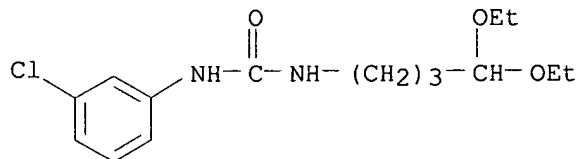


RN 306770-65-0 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-phenoxyphenyl)- (9CI) (CA INDEX NAME)



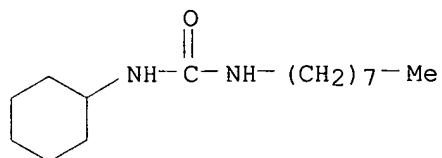
RN 306770-66-1 HCAPLUS

CN Urea, N-(3-chlorophenyl)-N'-(4,4-diethoxybutyl)- (9CI) (CA INDEX NAME)



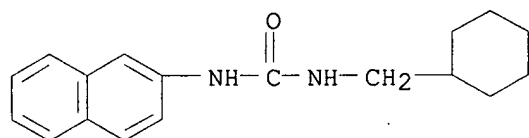
RN 306770-67-2 HCAPLUS

CN Urea, N-cyclohexyl-N'-octyl- (9CI) (CA INDEX NAME)



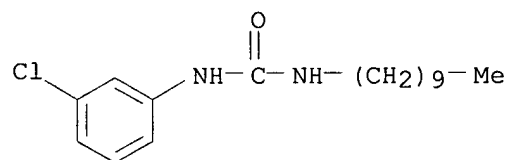
RN 306770-68-3 HCAPLUS

CN Urea, N-(cyclohexylmethyl)-N'-2-naphthalenyl- (9CI) (CA INDEX NAME)



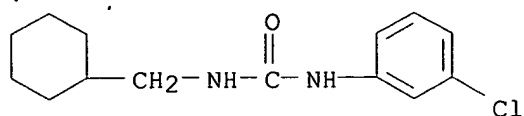
RN 306770-69-4 HCAPLUS

CN Urea, N-(3-chlorophenyl)-N'-decyl- (9CI) (CA INDEX NAME)

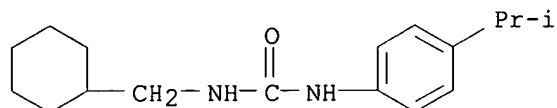


RN 306770-70-7 HCAPLUS

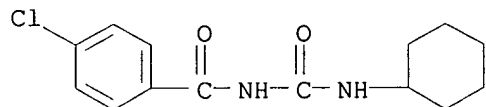
CN Urea, N-(3-chlorophenyl)-N'-(cyclohexylmethyl)- (9CI) (CA INDEX NAME)



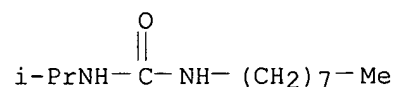
RN 306770-71-8 HCAPLUS  
 CN Urea, N-(cyclohexylmethyl)-N'-[4-(1-methylethyl)phenyl]- (9CI) (CA INDEX NAME)



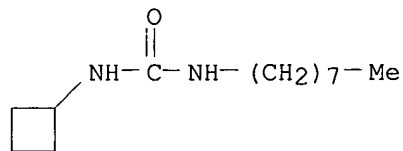
RN 326003-50-3 HCAPLUS  
 CN Benzamide, 4-chloro-N-[(cyclohexylamino)carbonyl]- (9CI) (CA INDEX NAME)



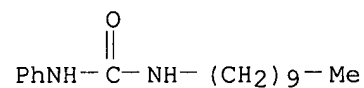
RN 338797-32-3 HCAPLUS  
 CN Urea, N-(1-methylethyl)-N'-octyl- (9CI) (CA INDEX NAME)



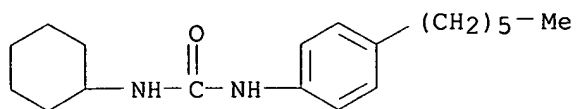
RN 338797-34-5 HCAPLUS  
 CN Urea, N-cyclobutyl-N'-octyl- (9CI) (CA INDEX NAME)



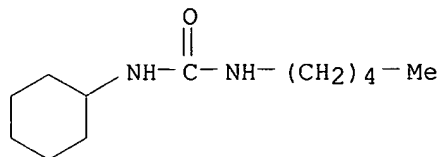
RN 339170-08-0 HCAPLUS  
 CN Urea, N-decyl-N'-phenyl- (9CI) (CA INDEX NAME)



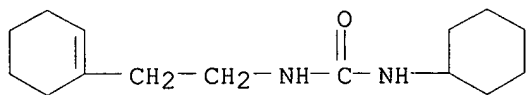
RN 392713-22-3 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-hexylphenyl)- (9CI) (CA INDEX NAME)



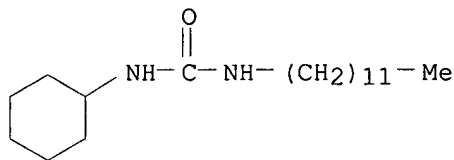
RN 400839-30-7 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-(4-pentyl- (9CI) (CA INDEX NAME)



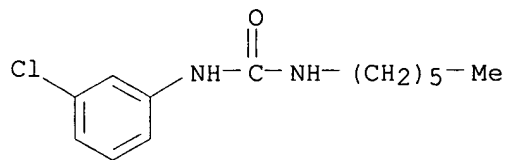
RN 401598-13-8 HCAPLUS  
 CN Urea, N-[2-(1-cyclohexen-1-yl)ethyl]-N'-cyclohexyl- (9CI) (CA INDEX NAME)



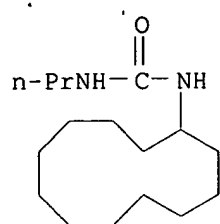
RN 402939-18-8 HCAPLUS  
 CN Urea, N-cyclohexyl-N'-dodecyl- (9CI) (CA INDEX NAME)



RN 446848-38-0 HCAPLUS  
 CN Urea, N-(3-chlorophenyl)-N'-hexyl- (9CI) (CA INDEX NAME)

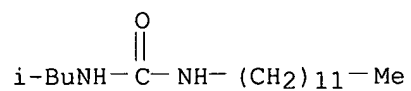


RN 447429-19-8 HCAPLUS  
 CN Urea, N-cyclododecyl-N'-propyl- (9CI) (CA INDEX NAME)



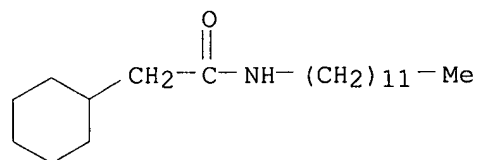
RN 460745-85-1 HCAPLUS

CN Urea, N-dodecyl-N'-(2-methylpropyl)- (9CI) (CA INDEX NAME)



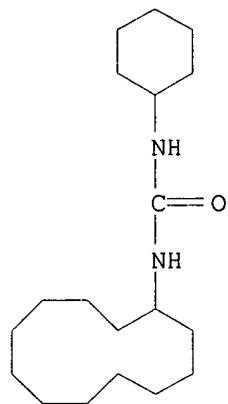
RN 479412-88-9 HCAPLUS

CN Cyclohexaneacetamide, N-dodecyl- (9CI) (CA INDEX NAME)



RN 479413-00-8 HCAPLUS

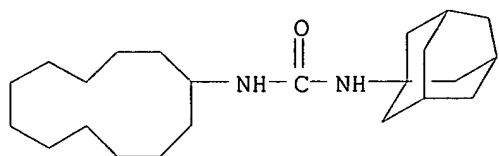
CN Urea, N-cyclododecyl-N'-cyclohexyl- (9CI) (CA INDEX NAME)



RN 479413-14-4 HCAPLUS

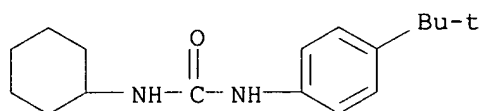
CN Urea, N-cyclododecyl-N'-tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl- (9CI) (CA INDEX NAME)





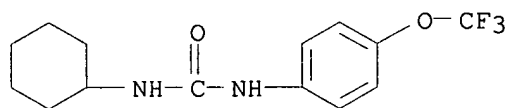
RN 479413-19-9 HCAPLUS

CN Urea, N-cyclohexyl-N'-[4-(1,1-dimethylethyl)phenyl]- (9CI) (CA INDEX NAME)



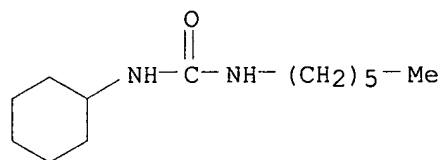
RN 479413-39-3 HCAPLUS

CN Urea, N-cyclohexyl-N'-[4-(trifluoromethoxy)phenyl]- (9CI) (CA INDEX NAME)



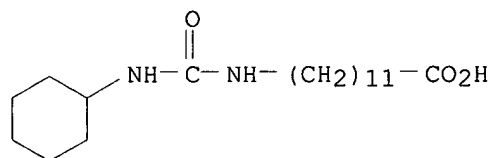
RN 479413-45-1 HCAPLUS

CN Urea, N-cyclohexyl-N'-hexyl- (9CI) (CA INDEX NAME)



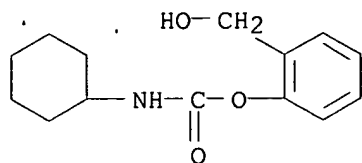
RN 479413-68-8 HCAPLUS

CN Dodecanoic acid, 12-[[cyclohexylamino]carbonylamino]- (9CI) (CA INDEX NAME)

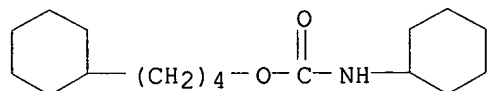


RN 501003-69-6 HCAPLUS

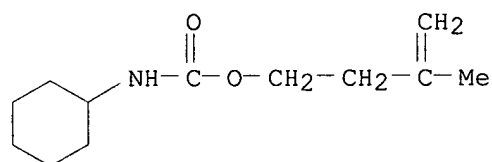
CN Carbamic acid, cyclohexyl-, 2-(hydroxymethyl)phenyl ester (9CI) (CA INDEX NAME)



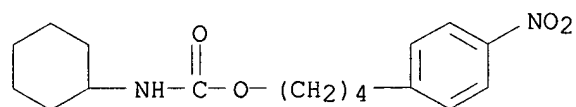
RN 501003-70-9 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, 4-cyclohexylbutyl ester (9CI) (CA INDEX NAME)



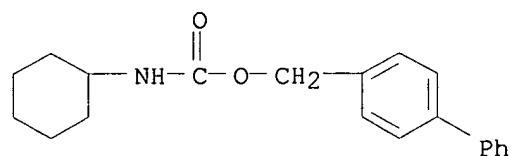
RN 501003-71-0 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, 3-methyl-3-butenyl ester (9CI) (CA INDEX NAME)



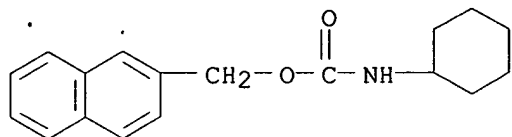
RN 501003-72-1 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, 4-(4-nitrophenyl)butyl ester (9CI) (CA INDEX NAME)



RN 501003-73-2 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, [1,1'-biphenyl]-4-ylmethyl ester (9CI) (CA INDEX NAME)

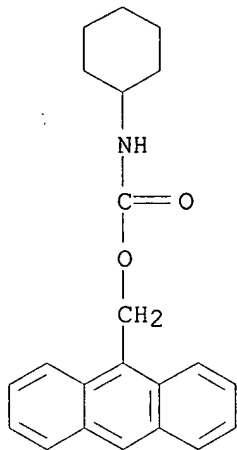


RN 501003-74-3 HCAPLUS  
 CN Carbamic acid, cyclohexyl-, 2-naphthalenylmethyl ester (9CI) (CA INDEX NAME)



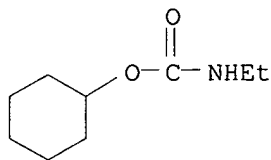
RN 501003-75-4 HCAPLUS

CN Carbamic acid, cyclohexyl-, 9-anthracenylmethyl ester (9CI) (CA INDEX NAME)

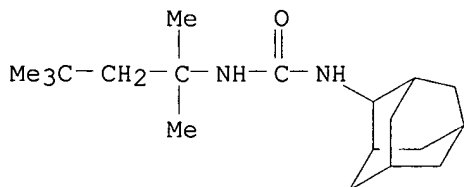


RN 501003-76-5 HCAPLUS

CN Carbamic acid, ethyl-, cyclohexyl ester (9CI) (CA INDEX NAME)

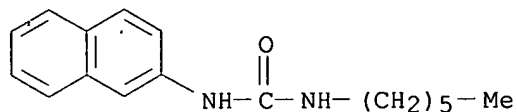


RN 501003-77-6 HCAPLUS

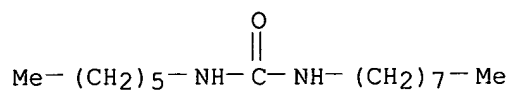
CN Urea, N-(1,1,3,3-tetramethylbutyl)-N'-tricyclo[3.3.1.1<sup>3,7</sup>]dec-2-yl- (9CI) (CA INDEX NAME)

RN 501003-78-7 HCAPLUS

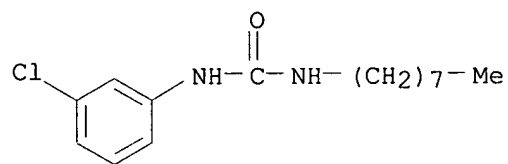
CN Urea, N-hexyl-N'-2-naphthalenyl- (9CI) (CA INDEX NAME)



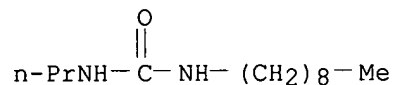
RN 501003-79-8 HCAPLUS  
 CN Urea, N-hexyl-N'-octyl- (9CI) (CA INDEX NAME)



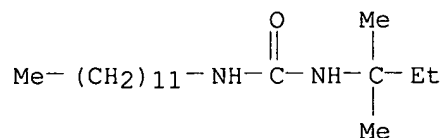
RN 501003-80-1 HCAPLUS  
 CN Urea, N-(3-chlorophenyl)-N'-octyl- (9CI) (CA INDEX NAME)



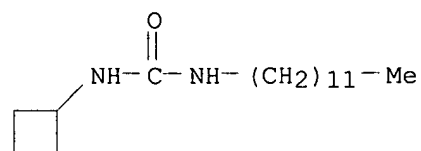
RN 501003-81-2 HCAPLUS  
 CN Urea, N-nonyl-N'-propyl- (9CI) (CA INDEX NAME)



RN 501003-82-3 HCAPLUS  
 CN Urea, N-(1,1-dimethylpropyl)-N'-dodecyl- (9CI) (CA INDEX NAME)

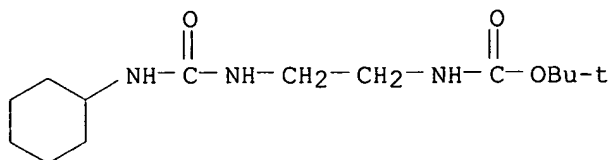


RN 501003-83-4 HCAPLUS  
 CN Urea, N-cyclobutyl-N'-dodecyl- (9CI) (CA INDEX NAME)



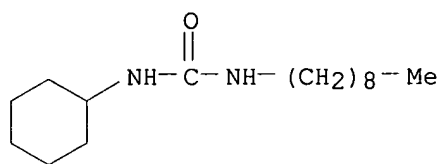
RN 501003-84-5 HCAPLUS

CN Carbamic acid, [2-[[[(cyclohexylamino)carbonyl]amino]ethyl]-,  
1,1-dimethylethyl ester (9CI) (CA INDEX NAME)



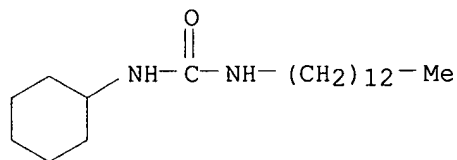
RN 501003-85-6 HCAPLUS

CN Urea, N-cyclohexyl-N'-nonyl- (9CI) (CA INDEX NAME)



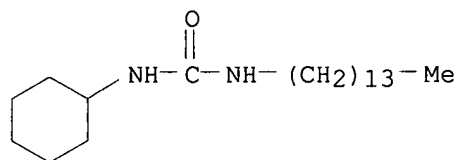
RN 501003-86-7 HCAPLUS

CN Urea, N-cyclohexyl-N'-tridecyl- (9CI) (CA INDEX NAME)



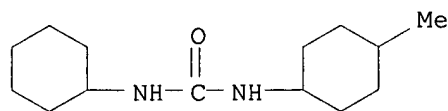
RN 501003-87-8 HCAPLUS

CN Urea, N-cyclohexyl-N'-tetradecyl- (9CI) (CA INDEX NAME)



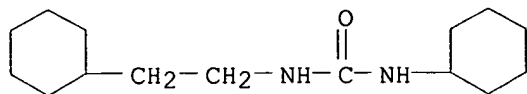
RN 501003-88-9 HCAPLUS

CN Urea, N-cyclohexyl-N'-(4-methylcyclohexyl)- (9CI) (CA INDEX NAME)



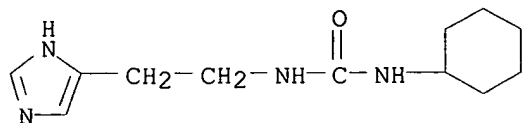
RN 501003-89-0 HCAPLUS

CN Urea, N-cyclohexyl-N'-(2-cyclohexylethyl)- (9CI) (CA INDEX NAME)



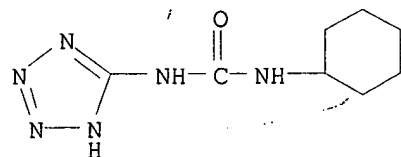
RN 501003-90-3 HCAPLUS

CN Urea, N-cyclohexyl-N'-[2-(1H-imidazol-4-yl)ethyl]- (9CI) (CA INDEX NAME)



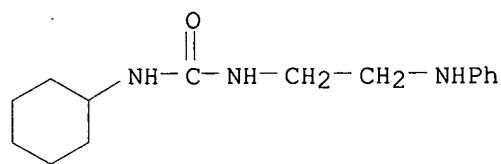
RN 501003-91-4 HCAPLUS

CN Urea, N-cyclohexyl-N'-1H-tetrazol-5-yl- (9CI) (CA INDEX NAME)



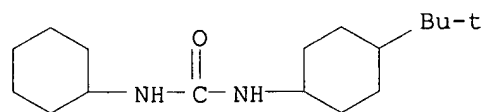
RN 501003-92-5 HCAPLUS

CN Urea, N-cyclohexyl-N'-[2-(phenylamino)ethyl]- (9CI) (CA INDEX NAME)



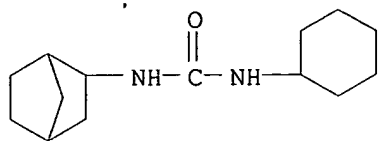
RN 501003-93-6 HCAPLUS

CN Urea, N-cyclohexyl-N'-[4-(1,1-dimethylethyl)cyclohexyl]- (9CI) (CA INDEX NAME)



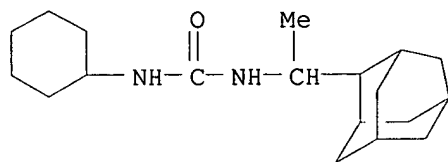
RN 501003-94-7 HCAPLUS

CN Urea, N-bicyclo[2.2.1]hept-2-yl-N'-cyclohexyl- (9CI) (CA INDEX NAME)



RN 501003-95-8 HCAPLUS

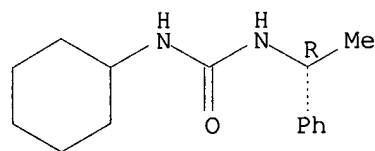
CN Urea, N'-cyclohexyl-N'-(1-tricyclo[3.3.1.1.3,7]dec-2-ylethyl)- (9CI) (CA INDEX NAME)



RN 501003-96-9 HCAPLUS

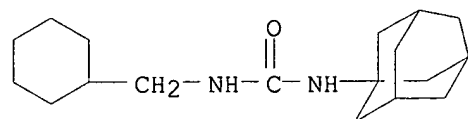
CN Urea, N-cyclohexyl-N'-[(1R)-1-phenylethyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).



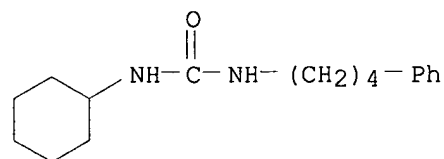
RN 501003-97-0 HCAPLUS

CN Urea, N-(cyclohexylmethyl)-N'-tricyclo[3.3.1.1.3,7]dec-1-yl- (9CI) (CA INDEX NAME)



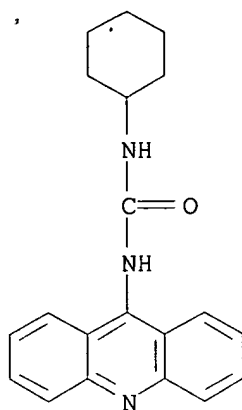
RN 501003-98-1 HCAPLUS

CN Urea, N-cyclohexyl-N'-(4-phenylbutyl)- (9CI) (CA INDEX NAME)



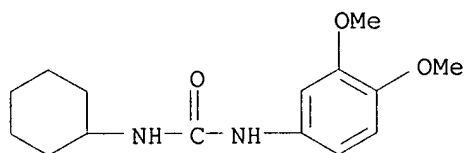
RN 501003-99-2 HCAPLUS

CN Urea, N-9-acridinyl-N'-cyclohexyl- (9CI) (CA INDEX NAME)



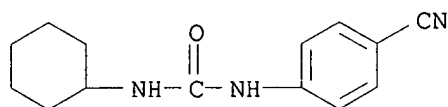
RN 501004-00-8 HCAPLUS

CN Urea, N-cyclohexyl-N'-(3,4-dimethoxyphenyl)- (9CI) (CA INDEX NAME)



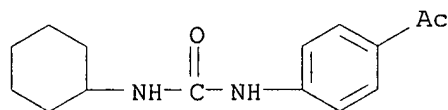
RN 501004-01-9 HCAPLUS

CN Urea, N-(4-cyanophenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



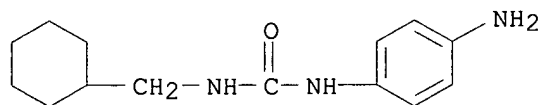
RN 501004-02-0 HCAPLUS

CN Urea, N-(4-acetylphenyl)-N'-cyclohexyl- (9CI) (CA INDEX NAME)



RN 501004-03-1 HCAPLUS

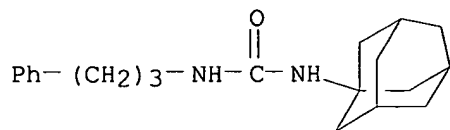
CN Urea, N-(4-aminophenyl)-N'-(cyclohexylmethyl)- (9CI) (CA INDEX NAME)



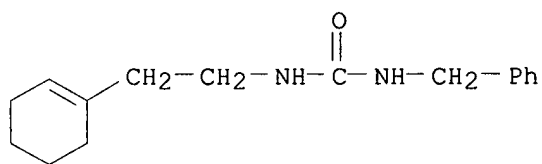
RN 501004-04-2 HCAPLUS



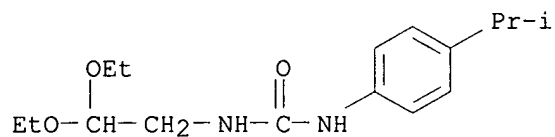
CN Urea, N-(3-phenylpropyl)-N'-tricyclo[3.3.1.1<sup>3,7</sup>]dec-1-yl- (9CI) (CA INDEX NAME)



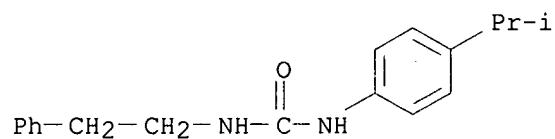
RN 501004-05-3 HCAPLUS  
CN Urea, N-[2-(1-cyclohexen-1-yl)ethyl]-N'-(phenylmethyl)- (9CI) (CA INDEX NAME)



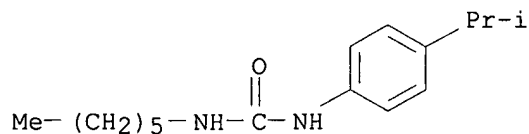
RN 501004-06-4 HCAPLUS  
CN Urea, N-(2,2-diethoxyethyl)-N'-[4-(1-methylethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 501004-07-5 HCAPLUS  
CN Urea, N-[4-(1-methylethyl)phenyl]-N'-(2-phenylethyl)- (9CI) (CA INDEX NAME)

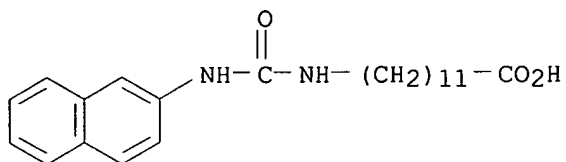


RN 501004-08-6 HCAPLUS  
CN Urea, N-hexyl-N'-[4-(1-methylethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 501004-09-7 HCAPLUS

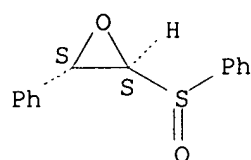
CN Dodecanoic acid, 12-[[ (2-naphthalenylamino)carbonyl]amino]- (9CI) (CA INDEX NAME)



RN 501004-10-0 HCAPLUS

CN Oxirane, 2-phenyl-3-(phenylsulfinyl)-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

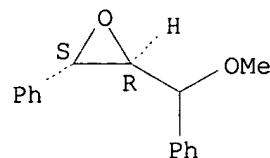
Relative stereochemistry.



RN 501004-11-1 HCAPLUS

CN Oxirane, 2-(methoxyphenylmethyl)-3-phenyl-, (2R,3S)-rel- (9CI) (CA INDEX NAME)

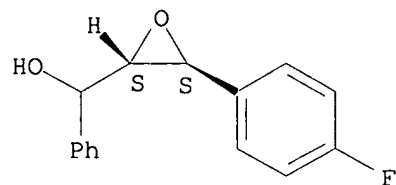
Relative stereochemistry.



RN 501004-12-2 HCAPLUS

CN Oxiranemethanol, 3-(4-fluorophenyl)-α-phenyl-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

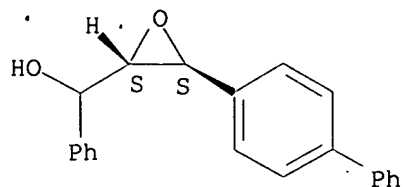
Relative stereochemistry.



RN 501004-13-3 HCAPLUS

CN Oxiranemethanol, 3-[1,1'-biphenyl]-4-yl-α-phenyl-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

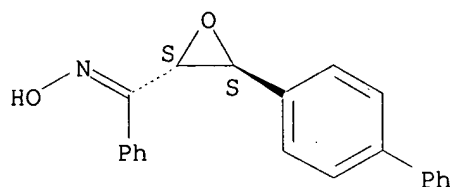
Relative stereochemistry.



RN 501004-14-4 HCAPLUS

CN Methanone, [(2R,3R)-3-[1,1'-biphenyl]-4-yloxiranyl]phenyl-, oxime, rel- (9CI) (CA INDEX NAME)

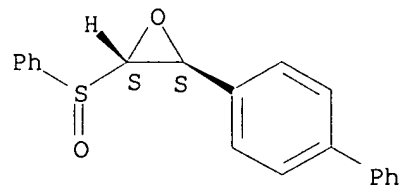
Relative stereochemistry.  
Double bond geometry unknown.



RN 501004-15-5 HCAPLUS

CN Oxirane, 2-[1,1'-biphenyl]-4-yl-3-(phenylsulfinyl)-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

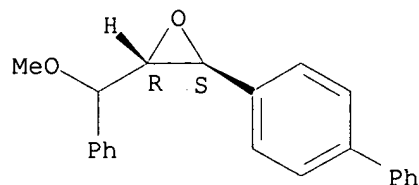
Relative stereochemistry.



RN 501004-16-6 HCAPLUS

CN Oxirane, 2-[1,1'-biphenyl]-4-yl-3-(methoxyphenylmethyl)-, (2R,3S)-rel- (9CI) (CA INDEX NAME)

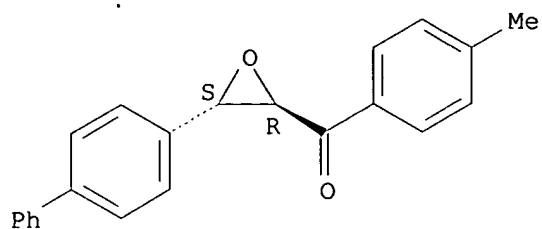
Relative stereochemistry.



RN 501004-17-7 HCAPLUS

CN Methanone, [(2R,3S)-3-[1,1'-biphenyl]-4-yloxiranyl](4-methylphenyl)-, rel- (9CI) (CA INDEX NAME)

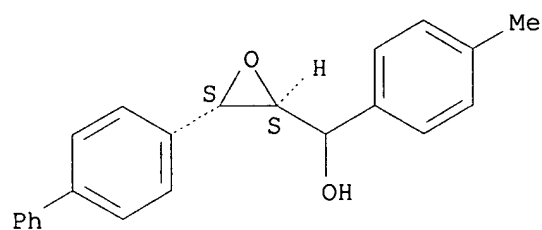
Relative stereochemistry.



RN 501004-18-8 HCAPLUS

CN Oxiranemethanol, 3-[1,1'-biphenyl]-4-yl- $\alpha$ -(4-methylphenyl)-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

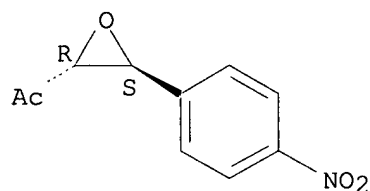
Relative stereochemistry.



RN 501004-19-9 HCAPLUS

CN Ethanone, 1-[(2R,3S)-3-(4-nitrophenyl)oxiranyl]-, rel- (9CI) (CA INDEX NAME)

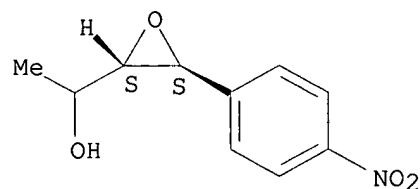
Relative stereochemistry.



RN 501004-20-2 HCAPLUS

CN Oxiranemethanol,  $\alpha$ -methyl-3-(4-nitrophenyl)-, (2R,3R)-rel- (9CI) (CA INDEX NAME)

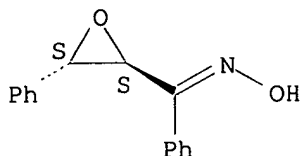
Relative stereochemistry.



RN 501011-71-8 HCAPLUS

CN Methanone, phenyl[(2R,3R)-3-phenyloxiranyl]-, oxime, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.  
Double bond geometry unknown.



REFERENCE COUNT: 43 THERE ARE 43 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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ACCESSION NUMBER: 2000:878331 HCAPLUS

DOCUMENT NUMBER: 134:54845

TITLE: Soluble epoxide hydrolase regulates hydrolysis of vasoactive epoxyeicosatrienoic acids

AUTHOR(S): Yu, Zhigang; Xu, Fengyun; Huse, Linn M.; **Morisseau, Christophe**; Draper, Alison J.; Newman, John W.; Parker, Carol; Graham, LeRae; Engler, Marguerite M.; **Hammock, Bruce D.**; **Zeldin, Darryl C.**; **Kroetz, Deanna L.**

CORPORATE SOURCE: Department of Biopharmaceutical Sciences, School of Nursing, University of California, San Francisco, CA, USA

SOURCE: Circulation Research (2000), 87(11), 992-998

CODEN: CIRUAL; ISSN: 0009-7330

PUBLISHER: Lippincott Williams & Wilkins

DOCUMENT TYPE: Journal

LANGUAGE: English

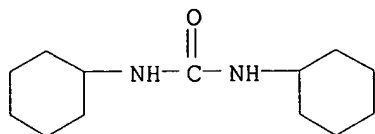
AB The cytochrome P 450-derived epoxyeicosatrienoic acids (EETs) have potent effects on renal vascular reactivity and tubular sodium and water transport; however, the role of these eicosanoids in the pathogenesis of hypertension is controversial. The current study examined the hydrolysis of the EETs to the corresponding dihydroxyeicosatrienoic acids (DHETs) as a mechanism for regulation of EET activity and blood pressure. EET hydrolysis was increased 5- to 54-fold in renal cortical S9 fractions from the spontaneously hypertensive rat (SHR) relative to the normotensive Wistar-Kyoto (WKY) rat. This increase was most significant for the 14,15-EET regioisomer, and there was a clear preference for hydrolysis of 14,15-EET over the 8,9- and 11,12-EETs. Increased EET hydrolysis was consistent with increased expression of soluble epoxide hydrolase (sEH) in the SHR renal microsomes and cytosol relative to the WKY samples. The urinary excretion of 14,15-DHET was 2.6-fold higher in the SHR than in the WKY rat, confirming increased EET hydrolysis in the SHR in vivo. Blood pressure was decreased 22 mm Hg 6 h after treatment of SHRs with the selective sEH inhibitor N,N'-dicyclohexylurea; this treatment had no effect on blood pressure in the WKY rat. These studies identify sEH as a novel therapeutic target for control of blood pressure. The identification of a potent and selective inhibitor of EET hydrolysis will be invaluable in separating the vascular effects of the EET and DHET eicosanoids.

IT **2387-23-7**, N,N'-Dicyclohexylurea

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

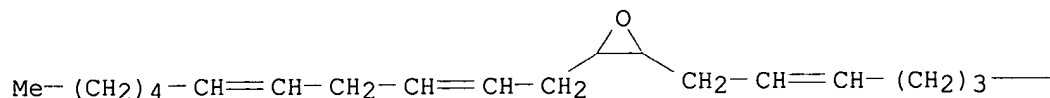
(soluble epoxide hydrolase inhibitor decrease of blood pressure in

hypertension)  
 RN 2387-23-7 HCAPLUS  
 CN Urea, N,N'-dicyclohexyl- (9CI) (CA INDEX NAME)



IT 81246-85-7 81276-02-0 81276-03-1  
 RL: ADV (Adverse effect, including toxicity); BPR (Biological process);  
 BSU (Biological study, unclassified); BIOL (Biological study); PROC  
 (Process)  
 (soluble epoxide hydrolase of kidney regulation of hydrolysis of  
 vasoactive epoxyeicosatrienoic acids in hypertension)  
 RN 81246-85-7 HCAPLUS  
 CN 5-Heptenoic acid, 7-[3-(2,5-undecadienyl)oxiranyl]- (9CI) (CA INDEX NAME)

PAGE 1-A

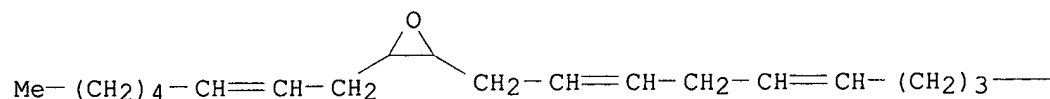


PAGE 1-B

—CO<sub>2</sub>H

RN 81276-02-0 HCAPLUS  
 CN 5,8-Decadienoic acid, 10-[3-(2-octenyl)oxiranyl]- (9CI) (CA INDEX NAME)

PAGE 1-A

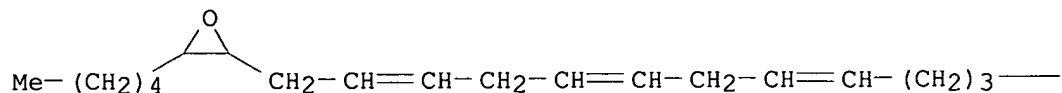


PAGE 1-B

—CO<sub>2</sub>H

RN 81276-03-1 HCAPLUS  
 CN 5,8,11-Tridecatricienoic acid, 13-(3-pentyloxiranyl)- (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

—CO<sub>2</sub>H

IT 79551-81-8 97717-67-4

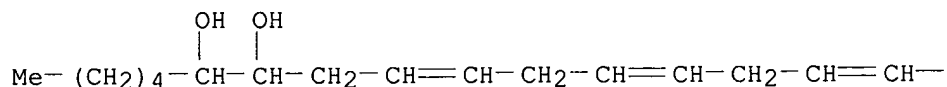
RL: ADV (Adverse effect, including toxicity); BSU (Biological study, unclassified); MFM (Metabolic formation); BIOL (Biological study); FORM (Formation, nonpreparative)

(soluble epoxide hydrolase of kidney regulation of hydrolysis of vasoactive epoxyeicosatrienoic acids in hypertension)

RN 79551-81-8 HCAPLUS

CN 5,8,11-Eicosatrienoic acid, 14,15-dihydroxy- (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

—(CH<sub>2</sub>)<sub>3</sub>-CO<sub>2</sub>H

RN 97717-67-4 HCAPLUS

CN Eicosatrienoic acid, dihydroxy- (9CI) (CA INDEX NAME)

CM 1

CRN 27613-79-2

CMF C20 H40 O4

CCI IDS

HO<sub>2</sub>C-(CH<sub>2</sub>)<sub>18</sub>-Me

2 ( D1-OH )

IT 9048-63-9, Epoxide hydrolase

RL: BAC (Biological activity or effector, except adverse); BOC (Biological occurrence); BPR (Biological process); BSU (Biological study,

unclassified); BIOL (Biological study); OCCU (Occurrence); PROC (Process)  
\* (soluble epoxide hydrolase of kidney regulation of hydrolysis of  
vasoactive epoxyeicosatrienoic acids in hypertension)  
RN 9048-63-9 HCAPLUS  
CN Hydratase, epoxide (9CI) (CA INDEX NAME)  
  
\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*  
IT 97717-69-6, Epoxyeicosatrienoic acid  
RL: ADV (Adverse effect, including toxicity); BPR (Biological process);  
BSU (Biological study, unclassified); BIOL (Biological study); PROC  
(Process)  
(triunsatd.; soluble epoxide hydrolase of kidney regulation of hydrolysis  
of vasoactive epoxyeicosatrienoic acids in hypertension)  
RN 97717-69-6 HCAPLUS  
CN Eicosatrienoic acid, epoxy- (9CI) (CA INDEX NAME)  
  
CM 1  
  
CRN 97717-68-5  
CMF C20 H38 O3  
CCI IDS

HO<sub>2</sub>C- (CH<sub>2</sub>)<sub>18</sub>-Me

D1-O-D1

REFERENCE COUNT: 29 THERE ARE 29 CITED REFERENCES AVAILABLE FOR THIS  
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT